AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions and listings of claims in the application:

LISTING OF CLAIMS:

1. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof, represented by formula [I]

$$X$$
 $COOR^1$
 $H_2N COOR^2$ [I]

Fwherein,

 R^1 and R^2 are identical or different, and each represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group substituted by one or two aryl groups, a hydroxy C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group), a group represented by formula- $CHR^cOC(O)ZR^d$ (wherein C_{1-10} alkyl group), a group represented by formula- $CHR^cOC(O)ZR^d$ (wherein C_{1-10} alkyl group, a C_{2-10} alkenyl group or an aryl group; and C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group, a group represented by formula [i]

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$$\begin{array}{c|c}
 & O \\
\hline
 & O \\
\hline
 & O \\
\hline
 & [i]
\end{array}$$

(wherein R^d is the same as described above) or a group represented by formula [ii]; or,

in the case where either R^1 or R^2 represents a hydrogen atom, the other represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{4-10} alkyl group substituted by one or two aryl groups, a hydroxy C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkoxy C_{1-10} alkyl group, a C_{1-10} alkyl group, a C_{1-10} alkyl group, a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are the same as described above), a group represented by formula- $CHR^cOC(O)ZR^d$ (wherein Z, R^c and R^d are the same as described above), a group represented by formula [i]

$$\begin{array}{c|c}
 & O \\
\hline
 &$$



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X represents a hydrogen atom or a fluorine atom; and

Y represents -OCHR³R⁴, -SR³, -S(O)_nR⁵, -SCHR³R⁴, -S(O)_nCHR³R⁴, -NHCHR³R⁴, -N(CHR³R⁴)(CHR³'R⁴), -NHCOR³ or -OCOR⁵ (wherein R³, R³', R⁴ and R⁴' are identical or different, and each represents a hydrogen atom, a C₁₋₁₀alkyl group, or a C₁₋₁₀alkenyl group, -a phenyl group, a naphthyl group, a naphthyl group substituted by one to seven halogen atoms, a heteroaromatic group or a phenyl group substituted by one to five substituents selected from a group consisting of a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group; R⁵ represents a C₁₋₁₀alkyl group, or a C₁₋₁₀alkenyl group, a phenyl group, a naphthyl group, a naphthyl group substituted by one to seven halogen atoms, a heteroaromatic group or a phenyl group substituted by one to five substituents selected from a group consisting of a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group; and n represents integer 1 or 2).

2. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof, represented by formula [II]

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Fwherein,

 R^1 and R^2 are identical or different, and each represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group substituted by one or two aryl groups, a hydroxy C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group), a group represented by formula- $CHR^cOC(O)ZR^d$ (wherein C_{1-10} alkyl group), a group represented by formula- C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group, and C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group, a group represented by formula [i]

$$\mathbb{R}^{d}$$
 O
 O
 O

in the case where either R^1 or R^2 represents a hydrogen atom, the other represents a C_{1-} ₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a C₁₋₁₀alkyl group substituted by one or two aryl groups, a hydroxyC₂₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, an aminoC₂₋₁₀alkyl group, a C₁₋₁₀alkoxyC₁₋₁₀alkyl group, a C₁₋₁₀alkoxycarbonylC₁₋₁₀alkyl group, a farnesyl group, a 4-morpholinylC₁₋₁₀alkyl group, a C₁₋₁₀alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^a-and R^b are the same as described above), or a group represented by formula-CHR OC(O)ZR (wherein Z, R and R are the same as described above), a group represented by formula [i]

$$R^{d}$$
 O
 O
 O

(wherein R^d is the same as described above) or a group represented by formula [ii];

X represents a hydrogen atom or a fluorine atom; and

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Y represents -OCHR³R⁴, -SR³, -S(O)_nR⁵, -SCHR³R⁴, -S(O)_nCHR³R⁴, -NHCHR³R⁴, -NHCHR³R⁴, -N(CHR³R⁴), -NHCOR³ or -OCOR⁵ (wherein R³, R³, R⁴ and R⁴ are identical or different, and each represents a hydrogen atom, a C₁₋₁₀alkyl group, <u>or</u> a C₁₋₁₀alkenyl group, <u>a</u> phenyl group, a naphthyl group, a naphthyl group substituted by one to seven halogen atoms, a heteroaromatic group or a phenyl group substituted by one to five substituents selected from a group consisting of a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group; R⁵ represents a C₁₋₁₀alkyl group, <u>or</u> a C₁₋₁₀alkenyl group, <u>a</u> cheteroaromatic group or a phenyl group substituted by one to seven halogen atoms, a heteroaromatic group or a phenyl group substituted by one to five substituents selected from a group consisting of a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group; and n represents integer 1 or 24.

3. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II],

 R^1 and R^2 are identical or different, and each represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group substituted by one or two phenyl groups, a hydroxy C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkoxy C_{1-10} alkyl group or a C_{1-10} alkoxycarbonyl C_{1-10} alkyl group; or,

in the case where either R^1 or R^2 represents a hydrogen atom, the other represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group substituted by one or two phenyl groups, a hydroxy C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl

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group, an amino C_{2-10} alkyl group, a C_{1-10} alkoxy C_{1-10} alkyl group or a - C_{1-10} alkoxycarbonyl C_{1-10} alkyl group.

4. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II],

 R^1 and R^2 are identical or different, and each represents a C_{1-10} alkyl group, a C_{2-6} alkenyl group, a C_{2-6} alkynyl group, a C_{1-6} alkyl group substituted by one or two phenyl groups, a hydroxy C_{2-6} alkyl group, a halogeno C_{1-6} alkyl group, an azido C_{1-6} alkyl group, an amino C_{2-6} alkyl group, a C_{1-6} alkoxy C_{1-6} alkyl group or a C_{1-6} alkoxycarbonyl C_{1-6} alkyl group; or,

in the case where either R^1 or R^2 represents a hydrogen atom, the other represents a C_{1-6} alkyl group, a C_{2-6} alkenyl group, a C_{2-6} alkynyl group, a C_{1-6} alkyl group substituted by one or two phenyl groups, a hydroxy C_{2-6} alkyl group, a halogeno C_{1-6} alkyl group, an azido C_{1-6} alkyl group, an amino C_{2-6} alkyl group, a C_{1-6} alkoxy C_{1-6} alkyl group or a C_{1-6} alkoxycarbonyl C_{1-6} alkyl group.

5. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II],

 R^1 and R^2 are identical or different, and each represents a farnesyl group, a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group), a group represented by formula- $C(O)ZR^d$ (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a

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 C_{1-10} alkyl group, <u>or</u> a C_{2-10} alkenyl group-or an aryl group; and R^d represents a C_{1-10} alkyl group, <u>or</u> a C_{2-10} alkenyl group-or an aryl group), a group represented by formula [i]

$$R^{d}$$
 O
 O
 O
 O

(wherein R^d is the same as described above) or a group represented by formula [ii]; or,

in the case where either R¹ or R² represents a hydrogen atom, the other represents a farnesyl group, a C₁₋₁₀alkyl group substituted by one or two aryl groups, a C₁₋₁₀alkyl group substituted by one or two aryl groups, a C₁₋₁₀alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^a and R^b are the same as described above), or a group represented by formula-CHR^cOC(O)ZR^d (wherein Z, R^c and R^d are the same as described above), a group represented by formula [i]

$$\begin{array}{c|c}
 & O \\
\hline
 &$$

6. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II],

 R^1 and R^2 are identical or different, and each represents a farnesyl group, a C_{1-6} alkyl group substituted by one or two aryl groups, a C_{1-6} alkoxycarbonyl C_{1-6} alkyl group, a 4morpholinyl C_{1-6} alkyl group, a C_{1-6} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-6} alkyl group), a group represented by formula-CHR $^cOC(O)ZR^d$ (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-6} alkyl group, or a C_{2-6} alkenyl group or an aryl group; and R^d represents a C_{1-6} alkyl group, or an aryl group), a group represented by formula [i]

$$\mathbb{R}^{d}$$
 O
 O
 O

in the case where either R¹ or R² represents a hydrogen atom, the other represents a farnesyl group, a C₁₋₆alkyl group substituted by one or two aryl groups, a C₁₋₆alkoxycarbonylC₁₋₆alkyl group, a 4-morpholinylC₁₋₆alkyl group, a C₁₋₁₀alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^a and R^b are the same as described above), or a group represented by formula-CHR^cOC(O)ZR^d (wherein Z,R^c and R^d are the same as described above), a group represented by formula [i]

$$\begin{array}{c|c}
 & O \\
\hline
 & O \\
\hline
 & O \\
\hline
 & [i]
\end{array}$$

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7. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom.

- 8. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; and X represents a fluorine atom.
- 9. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], wherein R² represents a hydrogen atom; and X represents a hydrogen atom.
- 10. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; and Y represents -OCHR³R⁴ (wherein R³ and R⁴ are the same as described above).
- 11. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; and Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above).
- **12. (currently amended):** A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, <u>or</u> a pharmaceutically acceptable salt thereof <u>or a hydrate thereof</u> according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; and Y represents -SR³ (wherein R³ is the same as described above).

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13. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; and Y represents -S(O)_nCHR³R⁴ (wherein R³, R⁴ and n are the same as described above).

- **14.** (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; and Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above).
- 15. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; and Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³, R⁴ and R⁴ are the same as described above).
- **16. (currently amended):** A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, <u>or</u> a pharmaceutically acceptable salt thereof <u>or a hydrate thereof</u> according to claim 2, wherein in the formula [II], represents a hydrogen atom; X represents a hydrogen atom; and Y represents -OCHR³R⁴ (wherein R³ and R⁴ are the same as described above).
- **17. (currently amended):** A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, <u>or</u> a pharmaceutically acceptable salt thereof <u>or a hydrate thereof</u> according to claim 2, wherein in the formula [II], represents a hydrogen atom; X represents a hydrogen atom; and Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above).
- **18.** (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim

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2, wherein in the formula [II], represents a hydrogen atom; X represents a hydrogen atom; and Y represents -SR³ (wherein R³ is the same as described above).

- **19.** (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; and Y represents -S(O)_nCHR³R⁴ (wherein R³, R⁴ and n are the same as described above).
- **20.** (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], wherein R² represents a hydrogen atom; X represents a hydrogen atom; and Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above).
- **21.** (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; and Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³', R⁴ and R⁴' are the same as described above).
- **22.** (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group, a C_{1-10} alkyl group, a farnesyl group, a C_{1-10} alkyl group or a $C_{$

 $_{10}$ alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

23. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a group represented by formula-CHR°OC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group), a group represented by formula [i]

$$\mathbb{R}^{d}$$
 O
 O
 O
 O

(wherein R^d is the same as described above) or a group represented by formula [ii]

24. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim

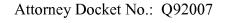
2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

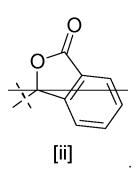
 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

25. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a group represented by formula-CHR^cOC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group), a group represented by formula [i]

$$\begin{array}{c|c}
 & O \\
\hline
 &$$





26. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents-SR³ (wherein R³ is the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

27. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -SR³ (wherein R³ is the same as described above); and

R¹ represents a group represented by formula-CHR^cOC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C₁₋₁₀alkyl group, <u>or</u> a C₂₋₁₀alkenyl group <u>or an aryl group</u>; and R^d represents a C₁₋₁₀alkyl group, <u>or</u> a C₂₋₁₀alkenyl group <u>or an aryl group</u>), a group represented by formula [i]

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$$\begin{array}{c|c}
 & O \\
\hline
 & O \\
\hline
 & O \\
\hline
 & [i]
\end{array}$$

(wherein R^d is the same as described above) or a group represented by formula [ii]

28. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -S(O)_nCHR³R⁴ (wherein R³, R⁴ and n are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

29. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim

2, wherein in the formula [II], R^2 represents a hydrogen atom; X represents a fluorine atom; Y represents $-S(O)_nCHR^3R^4$ (wherein R^3 , R^4 and n are the same as described above); and

 R^1 represents a group represented by formula-CHR^cOC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, <u>or</u> a C_{2-10} alkenyl group or an aryl group; and R^d represents a C_{1-10} alkyl group, <u>or</u> a C_{2-10} alkenyl group or an aryl group), a group represented by formula [i]

$$\mathbb{R}^{d}$$
 O
 O
 O
 O

(wherein R^d is the same as described above) or a group represented by formula [ii]

30. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group, a C

 $_{10}$ alkoxycarbonyl C_{1-10} alkyl group, a farnesyl group, a 4-morpholinyl C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

31. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a group represented by formula-CHR°OC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, <u>or</u> a C_{2-10} alkenyl group-or an aryl group; and R^d represents a C_{1-10} alkyl group, <u>or</u> a C_{2-10} alkenyl group-or an aryl group), a group represented by formula [i]

$$R^{d}$$
 O
 O
 O
 O

(wherein R^d is the same as described above) or a group represented by formula [ii]

32. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim

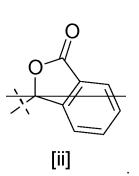
2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³, R⁴ and R⁴ are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

33. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³, R⁴ and R⁴ are the same as described above); and

 R^1 represents a group represented by formula-CHR COC(O)ZR (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group), a group represented by formula [i]

$$\mathbb{R}^{d}$$
 O
 O
 O



34. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents-OCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

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 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

35. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents-OCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

R¹ represents a group represented by formula-CHR^cOC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C₁₋₁₀alkyl group, <u>or a C₂₋₁₀alkenyl group or an aryl group</u>; and R^d represents a C₁₋₁₀alkyl group, <u>or a C₂₋₁₀alkenyl group or an aryl group</u>), a group represented by formula [i]

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$$R^{d}$$
 O
 O
 O
 O

(wherein R^d is the same as described above) or a group represented by formula [ii]

36. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group, a farnesyl group, a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

37. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim

2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a group represented by formula-CHR^cOC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond, R^c represents a hydrogen atom, C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or an aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or an aryl group), a group represented by formula [i]

$$\begin{array}{c|c}
 & O \\
\hline
 &$$

(wherein R^d is the same as described above) or a group represented by formula [ii]

38. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -SR³ (wherein R³ is the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group, a C

 $_{10}$ alkoxycarbonyl C_{1-10} alkyl group, a farnesyl group, a 4-morpholinyl C_{1-10} alkyl group, or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

39. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents-SR³ (wherein R³ is the same as described above); and

 R^1 represents a group represented by formula-CHR COC(O)ZR (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group), a group represented by formula [i]

$$\mathbb{R}^{d}$$
 \mathbb{Q}
 \mathbb{Q}

(wherein R^d is the same as described above) or a group represented by formula [ii]

40. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim

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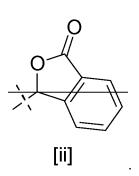
2, wherein in the formula [II], R^2 represents a hydrogen atom; X represents a hydrogen atom; Y represents $-S(O)_nCHR^3R^4$ (wherein R^3 , R^4 and n are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group substituted by one or two aryl groups, a hydroxy C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkoxy C_{1-10} alkyl group, a C_{1-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

41. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -S(O)_nCHR³R⁴ (wherein R³, R⁴ and n are the same as described above); and

R¹ represents a group represented by formula-CHR^cOC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom,C₁₋₁₀alkyl group, <u>or a C₂₋₁₀alkenyl group or an aryl group</u>; and R^d represents a C₁₋₁₀alkyl group, <u>or a</u> C₂₋₁₀alkenyl group or an aryl group), a group represented by formula [i]

$$R^{d}$$
 O
 O
 O
 O



42. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

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 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

43. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

R¹ represents a group represented by formula-CHR^cOC(O)XR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C₁₋₁₀alkyl group, <u>or</u> a C₂₋₁₀alkenyl group <u>or an aryl group</u>; and R^d represents a C₁₋₁₀alkyl group, <u>or</u> a C₂₋₁₀alkenyl group <u>or an aryl group</u>), a group represented by formula [i]

$$\begin{array}{c|c}
 & O \\
\hline
 & O \\
\hline
 & O \\
\hline
 & [i]
\end{array}$$

(wherein R^d is the same as described above) or a group represented by formula [ii]

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44. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³', R⁴ and R⁴' are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

45. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim

2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³', R⁴ and R⁴' are the same as described above); and

 R^1 represents a group represented by formula-CHR°OC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group; and R^d represents a C_{1-10} alkyl group, or an aryl group), a group represented by formula [i]

$$\mathbb{R}^{d}$$
 O
 O
 O

(wherein R^d is the same as described above) or a group represented by formula [ii]

46. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ represents a hydrogen atom; R⁴ represents a phenyl group or a phenyl group substituted by one to five substituents selected from a group coonsisting of a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a

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phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group substituted by one or two aryl groups, a hydroxy C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkoxy C_{1-10} alkyl group, a C_{1-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

47. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ represents a hydrogen atom; R⁴ represents a phenyl group or a phenyl group substituted by one to five substituents substituents selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

 R^1 represents a group represented by formula-CHR°OC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or an aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group), a group represented by formula [i]

$$\begin{array}{c|c}
 & O \\
\hline
 & O \\
\hline
 & O \\
\hline
 & [i]
\end{array}$$

(wherein R^d is the same as described above) or a group represented by formula [ii]

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- 48. (canceled).
- 49. (canceled).
- **50. (currently amended):** A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, <u>or</u> a pharmaceutically acceptable salt thereof <u>or a hydrate thereof</u> according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ and R⁴ are identical or different, and each represents a phenyl group or a phenyl group substituted by one to five substitutents selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

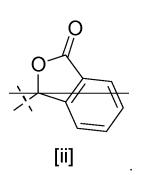
 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group, a C

 $_{10}$ alkoxycarbonyl C_{1-10} alkyl group, a farnesyl group, a 4-morpholinyl C_{1-10} alkyl group-or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

51. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, <u>or</u> a pharmaceutically acceptable salt thereof <u>or a hydrate thereof</u> according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ and R⁴ are identical or different, and each represents a phenyl group or a phenyl group substituted by one to five substitutents selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

 R^1 represents a group represented by formula-CHR^cOC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or an aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or an aryl group), a group represented by formula [i]

$$\begin{array}{c|c}
 & O \\
\hline
 & R^d & O
\end{array}$$
[i]



52. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, <u>or</u> a pharmaceutically acceptable salt thereof <u>or a hydrate thereof</u> according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -OCHR³R⁴ (wherein R³ represents a hydrogen atom; R⁴ represents a phenyl group or a phenyl group substituted by one to five substutuents selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, a maino group, a nitro group, a cyano group and a phenoxy group); and

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 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group a group group, a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

53. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -OCHR³R⁴ (wherein R³ represents a hydrogen atom; R⁴ represents a phenyl group or a

phenyl group substituted by one to five substutuents substituents selected from a group containing a halogen atom, a phenyl group, a C_{1-10} alkyl group, a C_{1-10} alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and phenoxy group); and

 R^1 represents a group represented by formula-CHR°OC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group), a group represented by formula [i]

$$\mathbb{R}^{d}$$
 O
 O
 O

- 54. (canceled).
- 55. (canceled).
- **56.** (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y

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represents -OCHR 3 R 4 (wherein R 3 and R 4 are identical or different, and each represents a phenyl group or a phenyl group substituted by one to five substituents selected from a group containing a halogen atom, a phenyl group, a C_{1-10} alkyl group, a C_{1-10} alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

57. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, <u>or</u> a pharmaceutically acceptable salt thereof <u>or a hydrate thereof</u> according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -OCHR³R⁴ (wherein R³ and R⁴ are identical or different, and each represents a phenyl group or a phenyl group substituted by one to five <u>substutuents substituents</u> selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, <u>a phenyl group</u>, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

 R^1 represents a group represented by formula-CHR°OC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, <u>or</u> a C_{2-10} alkenyl group or an aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group), a group represented by formula [i]

$$\begin{array}{c|c}
 & O \\
\hline
 & R^d & O
\end{array}$$

- **58.** (currently amended): A drug comprising the 2-amino-bicyclo [3.1.0] hexane 2,6-dicarboxylic ester-derivative, or the pharmaceutically acceptable salt thereof or the hydrate thereof according claim 2 as an active ingredient.
- **59. (original):** A drug according to claim 58, wherein the drag is a group II metabotropic glutamate receptor antagonist.
- **60. (previously presented):** (1R,2R,3R,5R,6R)-2-amino-3-(3,4-dichlorobenzyloxy)-6-fluoro-2,6-dicarboxylic acid 6-n-heptyl ester represented by the following structure: